

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110223\
 Data File : BN028433.D
 Acq On : 02 Nov 2023 22:10
 Operator : MA/JU
 Sample : 05196-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
ClientSampleId :
 PE-7-SOUTHSIDE-BASE

Manual Integrations
APPROVED
 Reviewed By :Yogesh Patel 11/04/2023
 Supervised By :mohammad ahmed 11/04/2023

Quant Time: Nov 03 01:50:40 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN110123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 02 06:26:50 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.724	152	12819	0.400	ng	0.00	
7) Naphthalene-d8	10.509	136	39414	0.400	ng	0.00	
13) Acenaphthene-d10	14.364	164	22372	0.400	ng	0.00	
19) Phenanthrene-d10	17.109	188	46607	0.400	ng	#-0.01	
29) Chrysene-d12	21.319	240	35946	0.400	ng	# 0.00	
35) Perylene-d12	23.631	264	31841	0.400	ng	#-0.01	
System Monitoring Compounds							
4) 2-Fluorophenol	5.305	112	9844	0.294	ng	0.00	
5) Phenol-d6	6.894	99	12594	0.297	ng	0.00	
8) Nitrobenzene-d5	8.875	82	7386	0.290	ng	0.00	
11) 2-Methylnaphthalene-d10	12.110	152	17267	0.306	ng	0.00	
14) 2,4,6-Tribromophenol	15.855	330	2264	0.260	ng	0.00	
15) 2-Fluorobiphenyl	13.006	172	4061	0.153	ng	0.00	
27) Fluoranthene-d10	19.152	212	34651	0.301	ng	0.00	
31) Terphenyl-d14	19.766	244	22764	0.301	ng	0.00	
Target Compounds							
9) Naphthalene	10.562	128	11624	0.109	ng	# 96	Qvalue
12) 2-Methylnaphthalene	12.182	142	6172	0.083	ng	# 93	
16) Acenaphthylene	14.086	152	25262	0.229	ng	98	
17) Acenaphthene	14.428	154	21430	0.306	ng	98	
18) Fluorene	15.411	166	18965	0.199	ng	# 97	
25) Phenanthrene	17.146	178	404117	2.935	ng	100	
26) Anthracene	17.245	178	84485	0.643	ng	98	
28) Fluoranthene	19.185	202	879391	5.472	ng	99	
30) Pyrene	19.547	202	799558	5.333	ng	# 95	
32) Benzo(a)anthracene	21.301	228	447276	3.135	ng	98	
33) Chrysene	21.355	228	414162	2.915	ng	98	
34) Bis(2-ethylhexyl)phtha...	21.266	149	186635	2.531	ng	99	
36) Indeno(1,2,3-cd)pyrene	26.002	276	217508	1.778	ng	# 94	
37) Benzo(b)fluoranthene	22.935	252	540566	4.423	ng	99	
38) Benzo(k)fluoranthene	22.979	252	184433m	1.479	ng		
39) Benzo(a)pyrene	23.529	252	341523	3.150	ng	99	
40) Dibenzo(a,h)anthracene	26.017	278	54783	0.538	ng	95	
41) Benzo(g,h,i)perylene	26.730	276	217449	1.994	ng	98	

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110223\
 Data File : BN028433.D
 Acq On : 02 Nov 2023 22:10
 Operator : MA/JU
 Sample : 05196-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PE-7-SOUTHSIDE-BASE

Quant Time: Nov 03 01:50:40 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN110123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 02 06:26:50 2023
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/04/2023
 Supervised By :mohammad ahmed 11/04/2023

