

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110222\
 Data File : BN022463.D
 Acq On : 02 Nov 2022 13:32
 Operator : CG/JU
 Sample : N5234-07DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 LSP-SS-06-20-40DL

Manual Integrations
 APPROVED

Reviewed By :Christian
 Giraldo

Quant Time: Nov 02 15:15:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN102722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 02 05:31:40 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							11/03/2022
1) 1,4-Dichlorobenzene-d4	7.928	152	3029	0.400	ng	0.00	Supervised By :mohammad
7) Naphthalene-d8	10.752	136	11096	0.400	ng	# 0.00	Ammed
13) Acenaphthene-d10	14.589	164	5422	0.400	ng	0.00	
19) Phenanthrene-d10	17.350	188	8844	0.400	ng	# 0.00	
29) Chrysene-d12	21.555	240	7629	0.400	ng	0.00	
35) Perylene-d12	23.999	264	7489	0.400	ng	0.00	11/07/2022
System Monitoring Compounds							
4) 2-Fluorophenol	5.458	112	641	0.079	ng	0.00	
5) Phenol-d6	7.105	99	1503	0.149	ng	0.01	
8) Nitrobenzene-d5	9.065	82	4858	0.506	ng	-0.04	
11) 2-Methylnaphthalene-d10	12.346	152	1007	0.058	ng	0.00	
14) 2,4,6-Tribromophenol	16.096	330	103	0.044	ng	0.00	
15) 2-Fluorobiphenyl	13.221	172	1037m	0.045	ng	0.00	
27) Fluoranthene-d10	19.390	212	907	0.037	ng	0.00	
31) Terphenyl-d14	19.985	244	959	0.048	ng	0.00	
Target Compounds							Qvalue
9) Naphthalene	10.795	128	40982	1.235	ng	# 80	
16) Acenaphthylene	14.311	152	19420	0.739	ng	# 78	
17) Acenaphthene	14.653	154	3277	0.183	ng	# 51	
18) Fluorene	15.648	166	12386	0.496	ng	# 87	
25) Phenanthrene	17.387	178	78451	2.587	ng	100	
26) Anthracene	17.487	178	23891	0.935	ng	97	
28) Fluoranthene	19.420	202	70995	2.159	ng	99	
30) Pyrene	19.782	202	80864	2.303	ng	97	
32) Benzo(a)anthracene	21.537	228	43686	1.455	ng	94	
33) Chrysene	21.591	228	38613	1.230	ng	96	
36) Indeno(1,2,3-cd)pyrene	26.561	276	23344	0.631	ng	95	
37) Benzo(b)fluoranthene	23.251	252	44966	1.312	ng	99	
38) Benzo(k)fluoranthene	23.292	252	15744m	0.462	ng		
39) Benzo(a)pyrene	23.888	252	42173m	1.512	ng		
40) Dibenzo(a,h)anthracene	26.572	278	6935	0.240	ng	# 75	
41) Benzo(g,h,i)perylene	27.359	276	23675	0.777	ng	98	

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

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