

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110122\
 Data File : BN022454.D
 Acq On : 02 Nov 2022 04:32
 Operator : CG/JU
 Sample : N5235-04
 Misc :
 ALS Vial : 62 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 LSP-DS-01-09-11

Manual Integrations
 APPROVED

Reviewed By :Christian
 Giraldo

Quant Time: Nov 02 05:41:20 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							11/02/2022
1) 1,4-Dichlorobenzene-d4	7.928	152	3905	0.400	ng	0.00	Supervised By :mohammad
7) Naphthalene-d8	10.752	136	11022	0.400	ng	0.00	Unmed
13) Acenaphthene-d10	14.589	164	4618	0.400	ng	0.00	
19) Phenanthrene-d10	17.350	188	7327	0.400	ng	# 0.00	
29) Chrysene-d12	21.546	240	8236	0.400	ng	# 0.00	
35) Perylene-d12	23.996	264	8387	0.400	ng	# 0.00	11/04/2022
System Monitoring Compounds							
4) 2-Fluorophenol	5.458	112	2730	0.262	ng	0.00	
5) Phenol-d6	7.090	99	3664	0.281	ng	0.00	
8) Nitrobenzene-d5	9.108	82	2217	0.233	ng	0.00	
11) 2-Methylnaphthalene-d10	12.346	152	3525	0.204	ng	0.00	
14) 2,4,6-Tribromophenol	16.096	330	452	0.228	ng	0.00	
15) 2-Fluorobiphenyl	13.221	172	4690	0.237	ng	0.00	
27) Fluoranthene-d10	19.386	212	5836	0.288	ng	0.00	
31) Terphenyl-d14	19.980	244	5950	0.277	ng	0.00	
Target Compounds							Qvalue
16) Acenaphthylene	14.311	152	7079	0.316	ng		98
18) Fluorene	15.647	166	882	0.042	ng	#	12
25) Phenanthrene	17.387	178	8113m	0.323	ng		
26) Anthracene	17.487	178	4606	0.218	ng		97
28) Fluoranthene	19.416	202	20758	0.762	ng		99
30) Pyrene	19.782	202	26033	0.687	ng	#	95
32) Benzo(a)anthracene	21.528	228	12637	0.390	ng	#	70
33) Chrysene	21.582	228	12614	0.372	ng	#	75
34) Bis(2-ethylhexyl)phtha...	21.448	149	14297	0.657	ng	#	87
36) Indeno(1,2,3-cd)pyrene	26.555	276	9353	0.226	ng		94
37) Benzo(b)fluoranthene	23.245	252	17205	0.448	ng	#	26
38) Benzo(k)fluoranthene	23.289	252	5706m	0.149	ng		
39) Benzo(a)pyrene	23.885	252	12668	0.406	ng	#	47
40) Dibenzo(a,h)anthracene	26.563	278	2885	0.089	ng	#	1
41) Benzo(g,h,i)perylene	27.353	276	13453	0.394	ng	#	75

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

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