

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122321\
 Data File : BN018183.D
 Acq On : 23 Dec 2021 16:01
 Operator : CG/JU
 Sample : M5179-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 POST-EXCAVATION-BOTTOM-WEST

Quant Time: Dec 24 23:31:51 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN122321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 23 14:22:48 2021
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 12/27/2021
 Supervised By :mohammad ahmed 12/30/2021

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.624	152	21728	0.400	ng	0.00
7) Naphthalene-d8	10.394	136	95867	0.400	ng	# 0.00
13) Acenaphthene-d10	14.256	164	60351	0.400	ng	0.00
19) Phenanthrene-d10	16.993	188	114404	0.400	ng	0.00
29) Chrysene-d12	21.196	240	105858	0.400	ng	0.00
35) Perylene-d12	23.420	264	102018	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.254	112	13335	0.281	ng	0.03
5) Phenol-d6	6.822	99	15289	0.301	ng	0.00
8) Nitrobenzene-d5	8.772	82	18069	0.280	ng	0.00
11) 2-Methylnaphthalene-d10	11.994	152	42155	0.256	ng	0.00
14) 2,4,6-Tribromophenol	15.753	330	5581	0.322	ng	0.00
15) 2-Fluorobiphenyl	12.886	172	64439	0.286	ng	0.00
27) Fluoranthene-d10	19.033	212	92326	0.255	ng	0.00
31) Terphenyl-d14	19.639	244	71925	0.309	ng	0.00
Target Compounds						
						Qvalue
9) Naphthalene	10.445	128	28119	0.119	ng	98
12) 2-Methylnaphthalene	12.070	142	12401	0.079	ng	99
16) Acenaphthylene	13.964	152	14902	0.066	ng	# 86
17) Acenaphthene	14.320	154	10087	0.065	ng	99
18) Fluorene	15.309	166	13889	0.072	ng	# 95
25) Phenanthrene	17.030	178	35259m	0.116	ng	
26) Anthracene	17.127	178	17844	0.071	ng	97
28) Fluoranthene	19.063	202	37997	0.107	ng	100
30) Pyrene	19.425	202	37258	0.110	ng	# 95
32) Benzo(a)anthracene	21.174	228	28194	0.092	ng	# 91
33) Chrysene	21.228	228	30097	0.090	ng	# 94
34) Bis(2-ethylhexyl)phtha...	21.121	149	589020	3.692	ng	99
36) Indeno(1,2,3-cd)pyrene	25.685	276	20182	0.068	ng	98
37) Benzo(b)fluoranthene	22.752	252	29572	0.094	ng	# 65
38) Benzo(k)fluoranthene	22.793	252	21979	0.069	ng	# 59
39) Benzo(a)pyrene	23.324	252	21504	0.076	ng	# 63
40) Dibenzo(a,h)anthracene	25.703	278	13106	0.082	ng	# 74
41) Benzo(g,h,i)perylene	26.377	276	21456	0.079	ng	# 90

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

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