

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111924\
 Data File : BN035204.D
 Acq On : 20 Nov 2024 18:07
 Operator : RC/JU
 Sample : PB165087BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS087

Quant Time: Nov 20 23:25:51 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN111624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 18 00:44:37 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.541	152	3122	0.400	ng/ul	0.00
4) Naphthalene-d8	10.300	136	7619	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.180	164	5223	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.937	188	11205	0.400	ng/ul	0.00
17) Chrysene-d12	21.148	240	7852	0.400	ng/ul	0.00
23) Perylene-d12	23.320	264	7902	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.136	96	1887	0.661	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.901	152	4509	0.404	ng/ul	-0.01
18) Fluoranthene-d10	18.977	212	10238	0.473	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.170	88	5392	1.741	ng/ul#	94
5) Naphthalene	10.350	128	7090	0.369	ng/ul	99
7) 2-Methylnaphthalene	11.978	142	5044	0.379	ng/ul	100
8) 1-Methylnaphthalene	12.198	142	5110	0.380	ng/ul	100
10) Acenaphthylene	13.898	152	7559	0.365	ng/ul	100
11) Acenaphthene	14.245	153	5907	0.373	ng/ul	99
12) Fluorene	15.240	166	6908	0.371	ng/ul	99
14) Pentachlorophenol	16.591	266	1500	0.575	ng/ul	98
15) Phenanthrene	16.975	178	11028	0.372	ng/ul	99
16) Anthracene	17.068	178	10002	0.364	ng/ul	100
19) Fluoranthene	19.005	202	12215	0.447	ng/ul	100
20) Pyrene	19.372	202	12311	0.433	ng/ul	100
21) Benzo(a)anthracene	21.131	228	10522	0.386	ng/ul	99
22) Chrysene	21.183	228	11079	0.404	ng/ul	99
24) Benzo(b)fluoranthene	22.674	252	10495	0.402	ng/ul	99
25) Benzo(k)fluoranthene	22.714	252	11026	0.409	ng/ul	100
26) Benzo(a)pyrene	23.226	252	9326	0.391	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.477	276	12688	0.394	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.490	278	9978	0.395	ng/ul	99
29) Benzo(g,h,i)perylene	26.131	276	10017	0.394	ng/ul	99

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

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