

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071224\
 Data File : BN032635.D
 Acq On : 11 Jul 2024 14:42
 Operator : MA/JU
 Sample : PB161917BS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS917

Quant Time: Jul 12 01:54:40 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN071224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 01:40:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.582	152	3963	0.400	ng/ul	0.00
4) Naphthalene-d8	10.365	136	13008	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.216	164	8501	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.986	188	18872	0.400	ng/ul	0.00
17) Chrysene-d12	21.197	240	15399	0.400	ng/ul	0.00
23) Perylene-d12	23.400	264	15446	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.106	96	3234	0.680	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.977	152	6557	0.334	ng/ul	0.00
18) Fluoranthene-d10	19.027	212	17120	0.341	ng/ul	0.00
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.139	88	9460	1.745	ng/ul	91
5) Naphthalene	10.415	128	12137	0.347	ng/ul	99
7) 2-Methylnaphthalene	12.059	142	7644	0.351	ng/ul	100
8) 1-Methylnaphthalene	12.252	142	7633	0.339	ng/ul	99
10) Acenaphthylene	13.943	152	11213	0.326	ng/ul	100
11) Acenaphthene	14.281	153	9951	0.350	ng/ul	99
12) Fluorene	15.294	166	10992	0.339	ng/ul	98
14) Pentachlorophenol	16.670	266	1944	0.587	ng/ul	99
15) Phenanthrene	17.033	178	17259	0.336	ng/ul	99
16) Anthracene	17.138	178	14113	0.336	ng/ul	100
19) Fluoranthene	19.055	202	21469	0.343	ng/ul	99
20) Pyrene	19.422	202	22545	0.348	ng/ul	99
21) Benzo(a)anthracene	21.188	228	16673	0.340	ng/ul	100
22) Chrysene	21.235	228	25124	0.354	ng/ul	100
24) Benzo(b)fluoranthene	22.737	252	18665	0.342	ng/ul	99
25) Benzo(k)fluoranthene	22.781	252	23556	0.355	ng/ul	99
26) Benzo(a)pyrene	23.310	252	18092	0.384	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.629	276	22262	0.349	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.639	278	17116	0.346	ng/ul	98
29) Benzo(g,h,i)perylene	26.304	276	18531	0.347	ng/ul	99

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

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