

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072122\
 Data File : BN020810.D
 Acq On : 21 Jul 2022 15:05
 Operator : CG/JU
 Sample : PB146317BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS317

Quant Time: Jul 21 15:34:36 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 20 01:02:30 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.825	152	2132	0.400	ng/ul	0.00
4) Naphthalene-d8	10.617	136	7268	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.469	164	5513	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.225	188	11244	0.400	ng/ul	0.00
17) Chrysene-d12	21.433	240	10233	0.400	ng/ul	0.00
23) Perylene-d12	23.786	264	8372	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.218	96	2092	0.847	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.223	152	4944	0.415	ng/ul	0.00
18) Fluoranthene-d10	19.262	212	11970	0.355	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.252	88	5127	2.033	ng/ul#	82
5) Naphthalene	10.667	128	7195	0.321	ng/ul	99
7) 2-Methylnaphthalene	12.294	142	5645	0.376	ng/ul	99
8) 1-Methylnaphthalene	12.509	142	6109	0.429	ng/ul	99
10) Acenaphthylene	14.192	152	8337	0.319	ng/ul	100
11) Acenaphthene	14.534	153	6846	0.327	ng/ul	100
12) Fluorene	15.529	166	7600	0.307	ng/ul	99
14) Pentachlorophenol	16.892	266	964	0.425	ng/ul	99
15) Phenanthrene	17.267	178	12925	0.331	ng/ul	100
16) Anthracene	17.360	178	10461	0.310	ng/ul	99
19) Fluoranthene	19.295	202	15116	0.323	ng/ul	100
20) Pyrene	19.657	202	15734	0.325	ng/ul	98
21) Benzo(a)anthracene	21.422	228	10761	0.293	ng/ul	100
22) Chrysene	21.471	228	14118	0.345	ng/ul	100
24) Benzo(b)fluoranthene	23.070	252	13279	0.343	ng/ul	99
25) Benzo(k)fluoranthene	23.117	252	13817	0.353	ng/ul	99
26) Benzo(a)pyrene	23.684	252	11799	0.335	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.220	276	13058	0.373	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.236	278	9992	0.353	ng/ul	98
29) Benzo(g,h,i)perylene	26.957	276	11739	0.390	ng/ul	99

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

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