

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062022\
 Data File : BN020285.D
 Acq On : 20 Jun 2022 19:05
 Operator : CG/JU
 Sample : PB145543BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS543

Quant Time: Jun 21 02:38:35 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN061822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 20 01:13:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.800	152	609	0.400	ng/ul	0.00
4) Naphthalene-d8	10.584	136	2474	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.423	164	2160	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.166	188	6032	0.400	ng/ul	0.00
17) Chrysene-d12	21.350	240	8769	0.400	ng/ul	0.00
23) Perylene-d12	23.662	264	10224	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.260	96	440	0.633	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.179	152	1388	0.339	ng/ul	0.00
18) Fluoranthene-d10	19.193	212	7132	0.394	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.294	88	1100	1.643	ng/ul#	76
5) Naphthalene	10.634	128	2296	0.290	ng/ul	97
7) 2-Methylnaphthalene	12.251	142	1594	0.311	ng/ul	100
8) 1-Methylnaphthalene	12.471	142	1789	0.355	ng/ul	99
10) Acenaphthylene	14.146	152	3147	0.305	ng/ul	98
11) Acenaphthene	14.488	153	2514	0.255	ng/ul	97
12) Fluorene	15.473	166	3204	0.313	ng/ul	99
14) Pentachlorophenol	16.833	266	760	0.656	ng/ul	99
15) Phenanthrene	17.209	178	6464	0.296	ng/ul	100
16) Anthracene	17.302	178	5557	0.301	ng/ul	99
19) Fluoranthene	19.225	202	9195	0.260	ng/ul	98
20) Pyrene	19.583	202	10078	0.263	ng/ul	100
21) Benzo(a)anthracene	21.336	228	10157	0.298	ng/ul	99
22) Chrysene	21.385	228	11629	0.274	ng/ul	100
24) Benzo(b)fluoranthene	22.960	252	14220	0.195	ng/ul	96
25) Benzo(k)fluoranthene	23.007	252	13997	0.206	ng/ul	97
26) Benzo(a)pyrene	23.560	252	13926	0.301	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.034	276	19202	0.255	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.047	278	15410	0.257	ng/ul	98
29) Benzo(g,h,i)perylene	26.755	276	17291	0.260	ng/ul	99

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

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