

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032524\
 Data File : BN030495.D
 Acq On : 25 Mar 2024 23:35
 Operator : MA/JU
 Sample : PB159747BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS747

Quant Time: Mar 26 00:32:53 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN032524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 25 13:38:40 2024
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.724	152	9361	0.400 ng/ul	0.00
4) Naphthalene-d8	10.507	136	26792	0.400 ng/ul	0.00
9) Acenaphthene-d10	14.363	164	15057	0.400 ng/ul	0.00
13) Phenanthrene-d10	17.108	188	34774	0.400 ng/ul	0.00
17) Chrysene-d12	21.301	240	24447	0.400 ng/ul	0.00
23) Perylene-d12	23.552	264	21605	0.400 ng/ul	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.218	96	6960	0.633 ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.102	152	12720	0.337 ng/ul	0.00
18) Fluoranthene-d10	19.137	212	27110	0.355 ng/ul	0.00
Target Compounds					
					Qvalue
2) 1,4-Dioxane	3.252	88	19087	1.679 ng/ul#	86
5) Naphthalene	10.557	128	25057	0.328 ng/ul	100
7) 2-Methylnaphthalene	12.174	142	16472	0.333 ng/ul	100
8) 1-Methylnaphthalene	12.399	142	16815	0.321 ng/ul	99
10) Acenaphthylene	14.081	152	24569	0.322 ng/ul	99
11) Acenaphthene	14.423	153	19121	0.325 ng/ul	99
12) Fluorene	15.413	166	21952	0.326 ng/ul	98
14) Pentachlorophenol	16.753	266	4236	0.647 ng/ul	99
15) Phenanthrene	17.150	178	37461	0.321 ng/ul	99
16) Anthracene	17.239	178	31500	0.319 ng/ul	100
19) Fluoranthene	19.170	202	40694	0.350 ng/ul	100
20) Pyrene	19.532	202	40251	0.351 ng/ul	99
21) Benzo(a)anthracene	21.283	228	33243	0.330 ng/ul	100
22) Chrysene	21.336	228	37039	0.347 ng/ul	100
24) Benzo(b)fluoranthene	22.870	252	33008	0.323 ng/ul	98
25) Benzo(k)fluoranthene	22.917	252	36556	0.350 ng/ul	98
26) Benzo(a)pyrene	23.449	252	27164	0.332 ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.826	276	32229	0.325 ng/ul	99
28) Dibenzo(a,h)anthracene	25.850	278	25951	0.327 ng/ul	98
29) Benzo(g,h,i)perylene	26.521	276	26243	0.306 ng/ul	99

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

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