

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN013123\
 Data File : BN023868.D
 Acq On : 31 Jan 2023 19:16
 Operator : CG/JU
 Sample : PB150492BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS492

Quant Time: Feb 01 01:04:18 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN013023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 01 01:00:14 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	3770	0.400	ng/ul	0.00
4) Naphthalene-d8	10.665	136	10361	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.510	164	6618	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.258	188	15272	0.400	ng/ul	0.00
17) Chrysene-d12	21.450	240	14912	0.400	ng/ul	0.00
23) Perylene-d12	23.855	264	13037	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.248	96	2288	0.566	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.260	152	4954	0.306	ng/ul	0.00
18) Fluoranthene-d10	19.284	212	13267	0.331	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.281	88	5933	1.386	ng/ul#	82
5) Naphthalene	10.715	128	8267	0.292	ng/ul	99
7) 2-Methylnaphthalene	12.332	142	5461	0.296	ng/ul	100
8) 1-Methylnaphthalene	12.552	142	5850	0.310	ng/ul	100
10) Acenaphthylene	14.232	152	7253	0.286	ng/ul	99
11) Acenaphthene	14.570	153	6513	0.292	ng/ul	99
12) Fluorene	15.560	166	7726	0.287	ng/ul	98
14) Pentachlorophenol	16.908	266	2462	0.489	ng/ul	99
15) Phenanthrene	17.300	178	13274	0.291	ng/ul	99
16) Anthracene	17.389	178	11396	0.299	ng/ul	99
19) Fluoranthene	19.317	202	16599	0.335	ng/ul	99
20) Pyrene	19.679	202	16832	0.328	ng/ul	99
21) Benzo(a)anthracene	21.435	228	16562	0.327	ng/ul	99
22) Chrysene	21.488	228	17312	0.325	ng/ul	99
24) Benzo(b)fluoranthene	23.124	252	17808	0.309	ng/ul	99
25) Benzo(k)fluoranthene	23.168	252	17834	0.317	ng/ul	99
26) Benzo(a)pyrene	23.741	252	14804	0.309	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.333	276	19778	0.294	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.349	278	16050	0.298	ng/ul	96
29) Benzo(g,h,i)perylene	27.101	276	16700	0.302	ng/ul	99

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

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