

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072624\
 Data File : BN033014.D
 Acq On : 27 Jul 2024 01:27
 Operator : MA/JU
 Sample : PB162101BS
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS101

Quant Time: Jul 27 02:05:13 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN072524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 25 15:43:00 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.537	152	2727	0.400	ng/ul	0.00
4) Naphthalene-d8	10.311	136	11397	0.400	ng/ul	0.01
9) Acenaphthene-d10	14.166	164	7492	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.945	188	15904	0.400	ng/ul	0.00
17) Chrysene-d12	21.154	240	14616	0.400	ng/ul	0.00
23) Perylene-d12	23.340	264	16348	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.069	96	2138	0.620	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.923	152	5077	0.273	ng/ul	0.01
18) Fluoranthene-d10	18.977	212	13049	0.315	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.098	88	6282	1.597	ng/ul#	70
5) Naphthalene	10.361	128	8888	0.304	ng/ul	99
7) 2-Methylnaphthalene	11.994	142	5579	0.281	ng/ul	99
8) 1-Methylnaphthalene	12.198	142	5975	0.279	ng/ul	100
10) Acenaphthylene	13.898	152	8949	0.292	ng/ul	99
11) Acenaphthene	14.231	153	7448	0.323	ng/ul	99
12) Fluorene	15.249	166	7891	0.294	ng/ul	100
14) Pentachlorophenol	16.633	266	1515	0.468	ng/ul	99
15) Phenanthrene	16.988	178	12196	0.286	ng/ul	99
16) Anthracene	17.093	178	10439	0.272	ng/ul	99
19) Fluoranthene	19.009	202	15875	0.330	ng/ul	99
20) Pyrene	19.376	202	17287	0.335	ng/ul	99
21) Benzo(a)anthracene	21.140	228	14650	0.294	ng/ul	99
22) Chrysene	21.192	228	18873	0.324	ng/ul	98
24) Benzo(b)fluoranthene	22.685	252	15668	0.328	ng/ul	96
25) Benzo(k)fluoranthene	22.729	252	19644	0.343	ng/ul	96
26) Benzo(a)pyrene	23.255	252	16062	0.357	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	25.554	276	18629	0.273	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.550	278	14470	0.267	ng/ul	97
29) Benzo(g,h,i)perylene	26.228	276	14709	0.271	ng/ul	98

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

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