

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090324\
 Data File : BN033745.D
 Acq On : 03 Sep 2024 17:34
 Operator : MA/JU
 Sample : PB163096BS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS096

Quant Time: Sep 04 01:42:22 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN082724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 04 01:39:37 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.523	152	6249	0.400	ng/ul	0.00
4) Naphthalene-d8	10.283	136	16977	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.161	164	8365	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.915	188	17148	0.400	ng/ul	0.00
17) Chrysene-d12	21.121	240	13639	0.400	ng/ul	0.00
23) Perylene-d12	23.275	264	14699	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.135	96	5215	0.812	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.878	152	9946	0.408	ng/ul	0.00
18) Fluoranthene-d10	18.952	212	18811	0.463	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.169	88	13612	1.910	ng/ul#	78
5) Naphthalene	10.327	128	16752	0.369	ng/ul	100
7) 2-Methylnaphthalene	11.955	142	10515	0.366	ng/ul	100
8) 1-Methylnaphthalene	12.175	142	12399	0.419	ng/ul	99
10) Acenaphthylene	13.874	152	13466	0.354	ng/ul	99
11) Acenaphthene	14.225	153	10817	0.394	ng/ul	99
12) Fluorene	15.215	166	11684	0.381	ng/ul	99
14) Pentachlorophenol	16.568	266	3255	0.603	ng/ul	99
15) Phenanthrene	16.953	178	18434	0.364	ng/ul	99
16) Anthracene	17.046	178	16591	0.365	ng/ul	99
19) Fluoranthene	18.980	202	20915	0.389	ng/ul	99
20) Pyrene	19.343	202	22072	0.386	ng/ul	98
21) Benzo(a)anthracene	21.103	228	20251	0.377	ng/ul	99
22) Chrysene	21.159	228	20691	0.381	ng/ul	99
24) Benzo(b)fluoranthene	22.637	252	20808	0.365	ng/ul	99
25) Benzo(k)fluoranthene	22.678	252	21478	0.376	ng/ul	98
26) Benzo(a)pyrene	23.181	252	19290	0.421	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.411	276	25198	0.366	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.425	278	19572	0.366	ng/ul	99
29) Benzo(g,h,i)perylene	26.059	276	19725	0.367	ng/ul	100

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

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