

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111624\
 Data File : BN035135.D
 Acq On : 16 Nov 2024 13:02
 Operator : RC/JU
 Sample : SSTD0.238
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.2224

Quant Time: Nov 18 00:34:46 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN111624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 18 00:33:27 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.545	152	2705	0.400	ng/ul	0.00
4) Naphthalene-d8	10.311	136	6167	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.189	164	3944	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.945	188	8580	0.400	ng/ul	0.00
17) Chrysene-d12	21.154	240	8005	0.400	ng/ul	0.00
23) Perylene-d12	23.331	264	8774	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.144	96	521	0.202	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.912	152	1798	0.190	ng/ul	0.00
18) Fluoranthene-d10	18.981	212	4335	0.211	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.174	88	575	0.204	ng/ul#	91
5) Naphthalene	10.361	128	3129	0.203	ng/ul	97
7) 2-Methylnaphthalene	11.983	142	2144	0.186	ng/ul	98
8) 1-Methylnaphthalene	12.203	142	2144	0.185	ng/ul	98
10) Acenaphthylene	13.903	152	3149	0.195	ng/ul	97
11) Acenaphthene	14.254	153	2415	0.198	ng/ul	99
12) Fluorene	15.244	166	2804	0.180	ng/ul	99
14) Pentachlorophenol	16.599	266	339	0.171	ng/ul	95
15) Phenanthrene	16.983	178	4559	0.199	ng/ul	97
16) Anthracene	17.076	178	4164	0.193	ng/ul	97
19) Fluoranthene	19.014	202	5446	0.204	ng/ul	99
20) Pyrene	19.376	202	5721	0.206	ng/ul	97
21) Benzo(a)anthracene	21.139	228	5550	0.191	ng/ul	98
22) Chrysene	21.189	228	5610	0.190	ng/ul	99
24) Benzo(b)fluoranthene	22.682	252	5733	0.186	ng/ul	91
25) Benzo(k)fluoranthene	22.726	252	6148	0.190	ng/ul#	91
26) Benzo(a)pyrene	23.235	252	5279	0.192	ng/ul#	88
27) Indeno(1,2,3-cd)pyrene	25.487	276	7135	0.212	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.507	278	5585	0.213	ng/ul	95
29) Benzo(g,h,i)perylene	26.147	276	5727	0.222	ng/ul	98

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

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