

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070524\
 Data File : BM046384.D
 Acq On : 05 Jul 2024 15:57
 Operator : MA/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV035

Quant Time: Jul 05 16:35:32 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM070524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 05 16:01:25 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.539	152	3485	0.400	ng/ul	0.00
4) Naphthalene-d8	10.304	136	9337	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.183	164	5685	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.934	188	12043	0.400	ng/ul	0.00
17) Chrysene-d12	21.140	240	8679	0.400	ng/ul	0.00
23) Perylene-d12	23.300	264	9162	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.130	96	1537	0.427	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.904	152	5525	0.400	ng/ul	0.00
18) Fluoranthene-d10	18.970	212	11614	0.382	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.164	88	1706	0.418	ng/ul	98
5) Naphthalene	10.354	128	10278	0.415	ng/ul	100
7) 2-Methylnaphthalene	11.976	142	7134	0.426	ng/ul	100
8) 1-Methylnaphthalene	12.201	142	6664	0.386	ng/ul	99
10) Acenaphthylene	13.896	152	11963	0.415	ng/ul	100
11) Acenaphthene	14.243	153	7672	0.408	ng/ul	99
12) Fluorene	15.238	166	9234	0.416	ng/ul	98
14) Pentachlorophenol	16.584	266	1314	0.374	ng/ul	99
15) Phenanthrene	16.977	178	14634	0.406	ng/ul	99
16) Anthracene	17.065	178	14718	0.421	ng/ul	99
19) Fluoranthene	19.002	202	16813	0.383	ng/ul	100
20) Pyrene	19.365	202	17788	0.394	ng/ul	96
21) Benzo(a)anthracene	21.125	228	15046	0.388	ng/ul	99
22) Chrysene	21.178	228	15257	0.396	ng/ul	100
24) Benzo(b)fluoranthene	22.662	252	14722	0.378	ng/ul	98
25) Benzo(k)fluoranthene	22.703	252	15229	0.387	ng/ul	99
26) Benzo(a)pyrene	23.206	252	13704	0.445	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.450	276	16147	0.388	ng/ul#	96
28) Dibenzo(a,h)anthracene	25.470	278	12656	0.393	ng/ul	97
29) Benzo(g,h,i)perylene	26.094	276	12828	0.387	ng/ul	97

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

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