

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120924\
 Data File : BM048913.D
 Acq On : 09 Dec 2024 17:55
 Operator : RC/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV012

Quant Time: Dec 09 21:43:36 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM120924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 09 17:52:26 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.586	152	3555	0.400	ng/ul	0.00
4) Naphthalene-d8	10.348	136	9131	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.220	164	4788	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.968	188	9764	0.400	ng/ul	0.00
17) Chrysene-d12	21.163	240	7760	0.400	ng/ul	0.00
23) Perylene-d12	23.335	264	8466	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	1730	0.416	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.948	152	4499	0.386	ng/ul	0.00
18) Fluoranthene-d10	19.002	212	8730	0.377	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.193	88	2173	0.452	ng/ul	95
5) Naphthalene	10.398	128	9158	0.391	ng/ul	99
7) 2-Methylnaphthalene	12.025	142	5618	0.393	ng/ul	100
8) 1-Methylnaphthalene	12.245	142	5368	0.369	ng/ul	99
10) Acenaphthylene	13.938	152	8190	0.405	ng/ul	99
11) Acenaphthene	14.285	153	6054	0.395	ng/ul	99
12) Fluorene	15.275	166	6565	0.386	ng/ul	96
14) Pentachlorophenol	16.626	266	675	0.336	ng/ul	94
15) Phenanthrene	17.010	178	10176	0.380	ng/ul	100
16) Anthracene	17.099	178	9053	0.379	ng/ul	99
19) Fluoranthene	19.030	202	12180	0.378	ng/ul	99
20) Pyrene	19.392	202	12909	0.372	ng/ul	98
21) Benzo(a)anthracene	21.146	228	9995	0.366	ng/ul	99
22) Chrysene	21.198	228	12451	0.403	ng/ul	100
24) Benzo(b)fluoranthene	22.689	252	11010	0.376	ng/ul	98
25) Benzo(k)fluoranthene	22.732	252	13403	0.396	ng/ul	98
26) Benzo(a)pyrene	23.238	252	10457	0.386	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.490	276	14762	0.388	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.507	278	11590	0.392	ng/ul	98
29) Benzo(g,h,i)perylene	26.138	276	11942	0.391	ng/ul	98

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

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