

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101824\
 Data File : BM048145.D
 Acq On : 18 Oct 2024 20:57
 Operator : RC/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV028

Quant Time: Oct 19 00:49:21 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM101824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 19 00:48:23 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.788	152	5936	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.585	136	18177	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.414	164	10613	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.162	188	18765	0.400	ng/ul	-0.01
17) Chrysene-d12	21.336	240	17594	0.400	ng/ul	0.00
23) Perylene-d12	23.595	264	17402	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	3270	0.405	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.201	152	9618	0.380	ng/ul	0.00
18) Fluoranthene-d10	19.183	212	20358	0.398	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.273	88	3607	0.381	ng/ul	99
5) Naphthalene	10.634	128	19840	0.402	ng/ul	99
7) 2-Methylnaphthalene	12.278	142	11628	0.411	ng/ul	99
8) 1-Methylnaphthalene	12.476	142	12059	0.371	ng/ul	99
10) Acenaphthylene	14.141	152	19911	0.406	ng/ul	98
11) Acenaphthene	14.479	153	13948	0.398	ng/ul	99
12) Fluorene	15.474	166	15297	0.408	ng/ul	99
14) Pentachlorophenol	16.837	266	2435	0.386	ng/ul	98
15) Phenanthrene	17.205	178	21031	0.443	ng/ul	100
16) Anthracene	17.306	178	20189	0.450	ng/ul	99
19) Fluoranthene	19.211	202	27373	0.396	ng/ul	99
20) Pyrene	19.574	202	29618	0.399	ng/ul	99
21) Benzo(a)anthracene	21.324	228	19934	0.368	ng/ul	99
22) Chrysene	21.368	228	31325	0.380	ng/ul	99
24) Benzo(b)fluoranthene	22.911	252	23759	0.356	ng/ul	96
25) Benzo(k)fluoranthene	22.955	252	30000	0.372	ng/ul	97
26) Benzo(a)pyrene	23.496	252	23640	0.426	ng/ul	93
27) Indeno(1,2,3-cd)pyrene	25.910	276	30560	0.367	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.923	278	23088	0.363	ng/ul	98
29) Benzo(g,h,i)perylene	26.607	276	25957	0.373	ng/ul	99

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

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