

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092524\
 Data File : BM047624.D
 Acq On : 25 Sep 2024 14:41
 Operator : RC/JU
 Sample : SST0.890
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.8018

Quant Time: Sep 25 15:24:35 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM092524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 25 15:01:59 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.830	152	5146	0.400	ng/ul	0.00
4) Naphthalene-d8	10.622	136	14273	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.459	164	7255	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.195	188	14600	0.400	ng/ul	0.00
17) Chrysene-d12	21.359	240	11950	0.400	ng/ul	0.00
23) Perylene-d12	23.639	264	12994	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.273	96	5400	0.823	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.211	152	15707	0.819	ng/ul	0.00
18) Fluoranthene-d10	19.215	212	30537	0.897	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.307	88	6421	0.821	ng/ul#	88
5) Naphthalene	10.671	128	31782	0.829	ng/ul	100
7) 2-Methylnaphthalene	12.288	142	19246	0.840	ng/ul	100
8) 1-Methylnaphthalene	12.502	142	20101	0.843	ng/ul	100
10) Acenaphthylene	14.177	152	28838	0.823	ng/ul	99
11) Acenaphthene	14.524	153	19887	0.819	ng/ul	99
12) Fluorene	15.505	166	22551	0.848	ng/ul	99
14) Pentachlorophenol	16.849	266	3484	0.837	ng/ul	99
15) Phenanthrene	17.237	178	35532	0.819	ng/ul	99
16) Anthracene	17.326	178	32251	0.835	ng/ul	100
19) Fluoranthene	19.247	202	42104	0.891	ng/ul	97
20) Pyrene	19.605	202	44329	0.890	ng/ul	99
21) Benzo(a)anthracene	21.342	228	40529	0.861	ng/ul	99
22) Chrysene	21.394	228	41385	0.823	ng/ul	100
24) Benzo(b)fluoranthene	22.949	252	43899	0.733	ng/ul	94
25) Benzo(k)fluoranthene	22.993	252	44521	0.774	ng/ul	94
26) Benzo(a)pyrene	23.540	252	36154	0.848	ng/ul	93
27) Indeno(1,2,3-cd)pyrene	25.960	276	55698	0.807	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.977	278	42515	0.808	ng/ul	97
29) Benzo(g,h,i)perylene	26.671	276	44514	0.799	ng/ul	99

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

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