

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031124\
 Data File : BM044574.D
 Acq On : 11 Mar 2024 12:19
 Operator : MA/JU
 Sample : SST0.812
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.8018

Quant Time: Mar 11 13:48:55 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM031124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 11 12:23:17 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.724	152	5758	0.400	ng/ul	0.00
4) Naphthalene-d8	10.501	136	18575	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.363	164	10661	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.120	188	23133	0.400	ng/ul	0.00
17) Chrysene-d12	21.312	240	16083	0.400	ng/ul	0.00
23) Perylene-d12	23.571	264	15419	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	6051	0.771	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.096	152	22136	0.858	ng/ul	0.00
18) Fluoranthene-d10	19.151	212	45070	0.911	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.314	88	6454	0.814	ng/ul#	76
5) Naphthalene	10.550	128	42744	0.796	ng/ul	99
7) 2-Methylnaphthalene	12.173	142	27634	0.829	ng/ul	98
8) 1-Methylnaphthalene	12.393	142	27726	0.823	ng/ul	100
10) Acenaphthylene	14.086	152	43688	0.770	ng/ul	100
11) Acenaphthene	14.428	153	31765	0.772	ng/ul	99
12) Fluorene	15.418	166	36603	0.787	ng/ul	99
14) Pentachlorophenol	16.770	266	4433	0.590	ng/ul	98
15) Phenanthrene	17.162	178	57610	0.805	ng/ul	99
16) Anthracene	17.255	178	54595	0.780	ng/ul	99
19) Fluoranthene	19.183	202	63454	0.857	ng/ul	99
20) Pyrene	19.546	202	64588	0.836	ng/ul	99
21) Benzo(a)anthracene	21.300	228	51981	0.742	ng/ul	98
22) Chrysene	21.350	228	54579	0.716	ng/ul	99
24) Benzo(b)fluoranthene	22.890	252	48991	0.796	ng/ul	94
25) Benzo(k)fluoranthene	22.937	252	52511	0.786	ng/ul	93
26) Benzo(a)pyrene	23.475	252	43629	0.769	ng/ul	91
27) Indeno(1,2,3-cd)pyrene	25.852	276	50666	0.711	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.879	278	39084	0.698	ng/ul	95
29) Benzo(g,h,i)perylene	26.547	276	43249	0.685	ng/ul	98

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

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