

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061521\
 Data File : BM030452.D
 Acq On : 15 Jun 2021 18:02
 Operator : CG/JU
 Sample : M2558-08
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X3554

Quant Time: Jun 15 18:33:16 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM060121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 15 11:43:56 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	4055	0.40	ng/ul	0.00
4) Naphthalene-d8	10.62	136	13880	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.45	164	7559	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.19	188	13901	0.40	ng/ul	0.00
17) Chrysene-d12	21.36	240	9905	0.40	ng/ul	0.00
23) Perylene-d12	23.64	264	9322	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.31	96	1858	0.47	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.20	152	5665	0.28	ng/ul	0.00
18) Fluoranthene-d10	19.21	212	12032	0.36	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	10.67	128	9252	0.225	ng/ul	98
7) 2-Methylnaphthalene	12.28	142	12254	0.442	ng/ul	99
10) Acenaphthylene	14.18	152	7516	0.209	ng/ul	98
11) Acenaphthene	14.52	153	8603	0.281	ng/ul	99
12) Fluorene	15.51	166	13545	0.389	ng/ul	99
14) Pentachlorophenol	16.86	266	2531	0.584	ng/ul	99
15) Phenanthrene	17.23	178	16548	0.329	ng/ul	99
16) Anthracene	17.33	178	17899	0.417	ng/ul	99
19) Fluoranthene	19.24	202	18682	0.370	ng/ul	97
20) Pyrene	19.61	202	18924	0.369	ng/ul	95
21) Benzo(a)anthracene	21.34	228	19130	0.481	ng/ul	100
22) Chrysene	21.40	228	19400	0.454	ng/ul	99
24) Benzo(b)fluoranthene	22.95	252	24784	0.535	ng/ul	96
25) Benzo(k)fluoranthene	22.99	252	12968	0.295	ng/ul	91
26) Benzo(a)pyrene	23.53	252	16298	0.436	ng/ul	97
28) Dibenzo(a,h)anthracene	25.96	278	26371	0.657	ng/ul	99
29) Benzo(g,h,i)perylene	26.65	276	12228	0.283	ng/ul	96

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

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