

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080521\
 Data File : BM031481.D
 Acq On : 05 Aug 2021 13:16
 Operator : CG/JU
 Sample : M3141-12
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGER0

Quant Time: Aug 05 13:47:28 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 05 11:03:39 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.590	152	1104	0.400	ng/ul	0.00
4) Naphthalene-d8	10.351	136	4430	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.219	164	2634	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.971	188	5099	0.400	ng/ul	0.00
17) Chrysene-d12	21.163	240	3953	0.400	ng/ul	0.00
23) Perylene-d12	23.321	264	3734	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.189	96	1979	1.614	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.947	152	1271	0.170	ng/ul	0.00
18) Fluoranthene-d10	19.004	212	2053	0.156	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.400	128	283	0.021	ng/ul#	52
10) Acenaphthylene	13.940	152	639	0.057	ng/ul#	81
15) Phenanthrene	17.012	178	920	0.057	ng/ul#	87
16) Anthracene	17.103	178	657	0.043	ng/ul#	76
19) Fluoranthene	19.031	202	2168	0.128	ng/ul#	89
21) Benzo(a)anthracene	21.146	228	1641	0.103	ng/ul#	85
22) Chrysene	21.197	228	1097	0.066	ng/ul#	85
24) Benzo(b)fluoranthene	22.684	252	1367	0.081	ng/ul#	24
25) Benzo(k)fluoranthene	22.723	252	639	0.037	ng/ul#	17

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

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