

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012622\
 Data File : BM033869.D
 Acq On : 26 Jan 2022 14:26
 Operator : CG/JU
 Sample : N1230-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0Q53

Quant Time: Jan 27 01:41:16 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM012422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 24 15:20:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	1820	0.400	ng/ul	0.00
4) Naphthalene-d8	10.752	136	6789	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.576	164	4198	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.313	188	8127	0.400	ng/ul	0.00
17) Chrysene-d12	21.475	240	6301	0.400	ng/ul #	0.00
23) Perylene-d12	23.871	264	6621	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	9365	4.922	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.339	152	3920	0.419	ng/ul	0.00
18) Fluoranthene-d10	19.331	212	9124	0.470	ng/ul	0.00
Target Compounds						
					Qvalue	
14) Pentachlorophenol	16.969	266	297	0.080	ng/ul	95
15) Phenanthrene	17.351	178	582	0.022	ng/ul#	75
19) Fluoranthene	19.361	202	643	0.022	ng/ul#	69
20) Pyrene	19.720	202	612	0.021	ng/ul#	21
21) Benzo(a)anthracene	21.458	228	652	0.025	ng/ul#	82
22) Chrysene	21.512	228	690	0.025	ng/ul#	82
24) Benzo(b)fluoranthene	23.139	252	787	0.026	ng/ul#	1
25) Benzo(k)fluoranthene	23.188	252	718	0.024	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	26.391	276	803	0.023	ng/ul#	85
28) Dibenzo(a,h)anthracene	26.405	278	571	0.021	ng/ul#	1
29) Benzo(g,h,i)perylene	27.154	276	762	0.026	ng/ul#	40

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

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