

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012622\
 Data File : BM033870.D
 Acq On : 26 Jan 2022 15:01
 Operator : CG/JU
 Sample : N1230-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0Q32

Quant Time: Jan 27 01:41:26 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	7.939	152	1845	0.400	ng/ul	0.00
4) Naphthalene-d8	10.751	136	6816	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.576	164	4039	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.313	188	7910	0.400	ng/ul	0.00
17) Chrysene-d12	21.477	240	6163	0.400	ng/ul #	0.00
23) Perylene-d12	23.873	264	6927	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.292	96	9475	4.912	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.344	152	3606	0.384	ng/ul	0.00
18) Fluoranthene-d10	19.334	212	8589	0.452	ng/ul	0.00
Target Compounds						
					Qvalue	
14) Pentachlorophenol	16.966	266	1685	0.464	ng/ul	97
15) Phenanthrene	17.354	178	813	0.032	ng/ul#	81
16) Anthracene	17.444	178	645	0.027	ng/ul#	78
19) Fluoranthene	19.361	202	1005	0.036	ng/ul#	77
20) Pyrene	19.723	202	884	0.031	ng/ul#	61
21) Benzo(a)anthracene	21.460	228	870	0.035	ng/ul#	79
22) Chrysene	21.514	228	896	0.034	ng/ul#	87
24) Benzo(b)fluoranthene	23.139	252	911	0.029	ng/ul#	1
25) Benzo(k)fluoranthene	23.187	252	826	0.026	ng/ul#	1
26) Benzo(a)pyrene	23.761	252	594	0.022	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	26.383	276	832	0.023	ng/ul#	90
28) Dibenzo(a,h)anthracene	26.402	278	683	0.024	ng/ul#	1
29) Benzo(g,h,i)perylene	27.144	276	713	0.023	ng/ul#	30

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

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