

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121522\
 Data File : BM038056.D
 Acq On : 16 Dec 2022 06:38
 Operator : CG/JU
 Sample : N5984-14
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 COLFO

Quant Time: Dec 16 07:11:30 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM121522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 15 21:53:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.059	152	5524	0.400	ng/ul	0.00
4) Naphthalene-d8	10.873	136	17327	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.685	164	13126	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.422	188	28500	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.585	240	19150	0.400	ng/ul	0.00
23) Perylene-d12	24.038	264	17563	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.406	96	21259	2.816	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.451	152	6987	0.276	ng/ul	0.00
18) Fluoranthene-d10	19.436	212	18373	0.265	ng/ul	0.00
Target Compounds						
					Qvalue	
7) 2-Methylnaphthalene	12.522	142	1757	0.054	ng/ul#	77
8) 1-Methylnaphthalene	12.742	142	878	0.027	ng/ul	98
14) Pentachlorophenol	17.088	266	4613932	391.003	ng/ul	99
15) Phenanthrene	17.464	178	2677	0.028	ng/ul#	1
16) Anthracene	17.552	178	7233	0.082	ng/ul#	21
19) Fluoranthene	19.464	202	4873	0.049	ng/ul#	42
20) Pyrene	19.831	202	79806	0.790	ng/ul#	95
21) Benzo(a)anthracene	21.565	228	2381	0.029	ng/ul#	1
22) Chrysene	21.600	228	7717	0.092	ng/ul#	72
24) Benzo(b)fluoranthene	23.281	252	3060	0.037	ng/ul#	20
25) Benzo(k)fluoranthene	23.333	252	2348	0.029	ng/ul#	1
26) Benzo(a)pyrene	23.927	252	1476	0.021	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	26.609	276	2157	0.024	ng/ul#	95
28) Dibenzo(a,h)anthracene	26.626	278	1636	0.024	ng/ul#	19
29) Benzo(g,h,i)perylene	27.400	276	1718	0.023	ng/ul#	62

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

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