

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080521\
 Data File : BM031581.D
 Acq On : 08 Aug 2021 10:30
 Operator : CG/JU
 Sample : M3163-02
 Misc :
 ALS Vial : 107 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGEX2

Quant Time: Aug 09 04:14:01 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 : Fri Aug 06 18:39:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.580	152	1177	0.400	ng/ul	0.00
4) Naphthalene-d8	10.337	136	4328	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.211	164	2452	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.960	188	4589	0.400	ng/ul	#-0.01
17) Chrysene-d12	21.157	240	3574	0.400	ng/ul	# 0.00
23) Perylene-d12	23.310	264	3788	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.186	96	4444	3.400	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.938	152	1694	0.232	ng/ul	0.00
18) Fluoranthene-d10	18.996	212	3147	0.264	ng/ul	0.00
Target Compounds						
					Qvalue	
14) Pentachlorophenol	16.620	266	39	0.027	ng/ul	96
15) Phenanthrene	17.002	178	343	0.024	ng/ul#	63
19) Fluoranthene	19.023	202	585	0.038	ng/ul#	64
20) Pyrene	19.389	202	575	0.037	ng/ul#	62
21) Benzo(a)anthracene	21.142	228	360	0.025	ng/ul#	68
22) Chrysene	21.196	228	390	0.026	ng/ul#	73
24) Benzo(b)fluoranthene	22.672	252	486	0.028	ng/ul#	1
26) Benzo(a)pyrene	23.213	252	430	0.029	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	25.456	276	409	0.021	ng/ul#	78
29) Benzo(g,h,i)perylene	26.101	276	355	0.022	ng/ul#	30

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

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