

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080521\
 Data File : BM031512.D
 Acq On : 06 Aug 2021 12:49
 Operator : CG/JU
 Sample : M3162-11
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGEW2

Manual Integrations
 APPROVED

mohammad
 8/6/2021 2:49:52 PM

Quant Time: Aug 06 13:20:39 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	7.587	152	1095	0.400	ng/ul	0.00
4) Naphthalene-d8	10.347	136	4038	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.216	164	2562	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.964	188	4777	0.400	ng/ul	#-0.01
17) Chrysene-d12	21.164	240	3989	0.400	ng/ul	0.00
23) Perylene-d12	23.319	264	4326	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.189	96	7861	6.464	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.939	152	3009	0.442	ng/ul	0.00
18) Fluoranthene-d10	19.001	212	5930	0.445	ng/ul	0.00
Target Compounds						
					Qvalue	
14) Pentachlorophenol	16.624	266	34	0.022	ng/ul#	86
15) Phenanthrene	17.009	178	448	0.030	ng/ul#	77
19) Fluoranthene	19.031	202	898	0.053	ng/ul#	76
20) Pyrene	19.393	202	807	0.047	ng/ul#	64
21) Benzo(a)anthracene	21.149	228	394	0.024	ng/ul#	67
22) Chrysene	21.197	228	463	0.028	ng/ul#	80
24) Benzo(b)fluoranthene	22.682	252	731m	0.038	ng/ul	
26) Benzo(a)pyrene	23.224	252	465	0.027	ng/ul#	1

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

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