

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080521\
 Data File : BM031513.D
 Acq On : 06 Aug 2021 13:26
 Operator : CG/JU
 Sample : M3162-20
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGF00

Manual Integrations
 APPROVED

mohammad
 8/6/2021 2:49:55 PM

Quant Time: Aug 06 14:00:53 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	1096	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	2464	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.968	188	4501	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.164	240	3445	0.400	ng/ul	0.00
23) Perylene-d12	23.319	264	3767	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.189	96	7649	6.284	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.944	152	2919	0.429	ng/ul	0.00
18) Fluoranthene-d10	19.001	212	5386	0.468	ng/ul	0.00
Target Compounds						
					Qvalue	
14) Pentachlorophenol	16.621	266	32	0.022	ng/ul#	71
19) Fluoranthene	19.027	202	695	0.047	ng/ul#	65
20) Pyrene	19.393	202	626	0.042	ng/ul#	57
21) Benzo(a)anthracene	21.147	228	442	0.032	ng/ul#	74
22) Chrysene	21.200	228	472	0.033	ng/ul#	74
24) Benzo(b)fluoranthene	22.679	252	601m	0.035	ng/ul	
26) Benzo(a)pyrene	23.224	252	411	0.028	ng/ul#	1

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

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