

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070821\
 Data File : BN015406.D
 Acq On : 08 Jul 2021 13:31
 Operator : CG/JU
 Sample : M2825-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGDQ6

Manual Integrations
 APPROVED

mohammad
 7/9/2021 3:03:30 PM

Quant Time: Jul 08 14:10:11 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 08 11:17:31 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.707	152	3150	0.400	ng/ul	0.00
4) Naphthalene-d8	10.480	136	10800	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.340	164	5848	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.086	188	12146m	0.400	ng/ul	0.00
17) Chrysene-d12	21.293	240	9391	0.400	ng/ul	0.00
23) Perylene-d12	23.534	264	9159	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.277	96	13958	3.931	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.069	152	5538	0.327	ng/ul	0.00
18) Fluoranthene-d10	19.122	212	11536	0.374	ng/ul	0.00
Target Compounds						
					Qvalue	
14) Pentachlorophenol	16.744	266	148	0.055	ng/ul	99
15) Phenanthrene	17.124	178	2122	0.048	ng/ul#	92
19) Fluoranthene	19.150	202	2530	0.058	ng/ul#	89
20) Pyrene	19.517	202	2250	0.052	ng/ul#	88
21) Benzo(a)anthracene	21.275	228	994	0.029	ng/ul#	87
22) Chrysene	21.328	228	960	0.023	ng/ul#	88
24) Benzo(b)fluoranthene	22.868	252	1369m	0.032	ng/ul	
26) Benzo(a)pyrene	23.438	252	922	0.026	ng/ul#	1

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

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