

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112221\
 Data File : BM033216.D
 Acq On : 22 Nov 2021 14:34
 Operator : CG/JU
 Sample : M4642-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X3805

Quant Time: Nov 22 15:05:24 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM111921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 19 15:41:12 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.956	152	2420	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.751	136	6676	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.576	164	3793	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.313	188	8433	0.400	ng/ul	0.00
17) Chrysene-d12	21.468	240	7744	0.400	ng/ul	0.00
23) Perylene-d12	23.810	264	7240	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.413	96	1471	0.646	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.334	152	3502	0.379	ng/ul	0.00
18) Fluoranthene-d10	19.327	212	8936	0.398	ng/ul	0.00
Target Compounds						
						Qvalue
5) Naphthalene	10.801	128	15479	0.748	ng/ul	98
7) 2-Methylnaphthalene	12.406	142	6683	0.499	ng/ul	99
12) Fluorene	15.619	166	11840	0.721	ng/ul	99
14) Pentachlorophenol	16.959	266	7631	2.711	ng/ul	98
16) Anthracene	17.445	178	24838	1.005	ng/ul	99
19) Fluoranthene	19.357	202	26224	0.704	ng/ul	98
20) Pyrene	19.720	202	30410	0.822	ng/ul	97
22) Chrysene	21.504	228	19785	0.651	ng/ul	99
24) Benzo(b)fluoranthene	23.098	252	18068	0.551	ng/ul	96
26) Benzo(a)pyrene	23.704	252	22379	0.824	ng/ul#	88
28) Dibenzo(a,h)anthracene	26.219	278	42056	1.537	ng/ul#	88

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

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