

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121721\
 Data File : BM033512.D
 Acq On : 17 Dec 2021 18:53
 Operator : CG/JU
 Sample : M5051-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P-3

Quant Time: Dec 18 04:54:58 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM121621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 16 17:10:15 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.937	152	46167	20.000	ng	0.00
21) Naphthalene-d8	10.748	136	187650	20.000	ng	# 0.00
39) Acenaphthene-d10	14.577	164	126313	20.000	ng	-0.01
64) Phenanthrene-d10	17.330	188	255706	20.000	ng	# 0.00
76) Chrysene-d12	21.506	240	157387	20.000	ng	0.00
86) Perylene-d12	23.918	264	132134	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.501	112	400541	149.706	ng	0.00
7) Phenol-d6	7.107	99	503108	140.121	ng	0.00
23) Nitrobenzene-d5	9.107	82	449179	102.780	ng	0.00
42) 2,4,6-Tribromophenol	16.071	330	208103	146.460	ng	0.00
45) 2-Fluorobiphenyl	13.207	172	913199	100.008	ng	0.00
79) Terphenyl-d14	19.953	244	1275025	127.540	ng	0.00
Target Compounds						
87) Indeno(1,2,3-cd)pyrene	26.447	276	15225	2.499	ng	# 86
91) Dibenzo(a,h)anthracene	26.441	278	12794	2.800	ng	97

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

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