

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121721\
 Data File : BM033507.D
 Acq On : 17 Dec 2021 15:52
 Operator : CG/JU
 Sample : M5064-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SASP2-121-1-(17-32)

Quant Time: Dec 18 04:51:48 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	51403	20.000	ng	0.00
21) Naphthalene-d8	10.748	136	210891	20.000	ng	0.00
39) Acenaphthene-d10	14.583	164	137267	20.000	ng	0.00
64) Phenanthrene-d10	17.330	188	267064	20.000	ng	0.00
76) Chrysene-d12	21.506	240	155390	20.000	ng	0.00
86) Perylene-d12	23.924	264	136044	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.502	112	353398	118.632	ng	0.00
7) Phenol-d6	7.107	99	486996	121.818	ng	0.00
23) Nitrobenzene-d5	9.107	82	377557	76.871	ng	0.00
42) 2,4,6-Tribromophenol	16.071	330	165373	107.100	ng	0.00
45) 2-Fluorobiphenyl	13.207	172	749755	75.556	ng	0.00
79) Terphenyl-d14	19.948	244	830488	84.141	ng	0.00
Target Compounds						
10) Phenol	7.131	94	12679	3.139	ng	Qvalue # 62

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

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