

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111821\
 Data File : BP008014.D
 Acq On : 18 Nov 2021 13:42
 Operator : CG/JU
 Sample : M4724-05
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 STUCCO-1ST-FLOOR

Quant Time: Nov 18 15:16:40 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 15 05:36:49 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	103593	20.000	ng	0.00
21) Naphthalene-d8	10.628	136	377338	20.000	ng	-0.01
39) Acenaphthene-d10	14.475	164	253668	20.000	ng	0.00
64) Phenanthrene-d10	17.228	188	574465	20.000	ng	#-0.01
76) Chrysene-d12	21.322	240	735150	20.000	ng	-0.01
86) Perylene-d12	23.722	264	845409	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.423	112	504969	79.076	ng	0.00
7) Phenol-d6	7.011	99	608571	71.635	ng	-0.01
23) Nitrobenzene-d5	8.993	82	387905	55.012	ng	-0.01
42) 2,4,6-Tribromophenol	15.969	330	335597	84.085	ng	0.00
45) 2-Fluorobiphenyl	13.099	172	968012	53.529	ng	0.00
79) Terphenyl-d14	19.792	244	1933741	53.337	ng	-0.01
Target Compounds						
60) Diethylphthalate	15.298	149	39641	2.049	ng	99
74) Di-n-butylphthalate	18.192	149	106252	3.178	ng	99

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

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