

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111821\
 Data File : BP008026.D
 Acq On : 18 Nov 2021 20:43
 Operator : CG/JU
 Sample : M4724-09
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 2ND-FLOOR-STUCCO

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 11/19/2021
 Supervised By :Sohil Jodhani 11/30/2021

Quant Time: Nov 19 02:25:01 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	112348	20.000	ng	0.00
21) Naphthalene-d8	10.628	136	389013	20.000	ng	#-0.01
39) Acenaphthene-d10	14.469	164	243912	20.000	ng	-0.01
64) Phenanthrene-d10	17.228	188	557150	20.000	ng	#-0.01
76) Chrysene-d12	21.322	240	705493	20.000	ng	#-0.01
86) Perylene-d12	23.716	264	813228	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	446006	64.400	ng	0.00
7) Phenol-d6	7.005	99	539519	58.558	ng	-0.02
23) Nitrobenzene-d5	8.987	82	338863	46.615	ng	-0.02
42) 2,4,6-Tribromophenol	15.963	330	279736	72.892	ng	-0.01
45) 2-Fluorobiphenyl	13.093	172	795331	45.739	ng	-0.01
79) Terphenyl-d14	19.792	244	1592094	45.760	ng	-0.01
Target Compounds						Qvalue
56) 4-Nitrophenol	14.716	139	10058m	3.019	ng	
74) Di-n-butylphthalate	18.192	149	78213	2.412	ng	98

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

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