

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111821\
 Data File : BP008032.D
 Acq On : 19 Nov 2021 00:06
 Operator : CG/JU
 Sample : M4724-13
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 3RD-FLOOR-BRICK

Quant Time: Nov 19 02:26:12 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	7.828	152	114514	20.000	ng	-0.01
21) Naphthalene-d8	10.628	136	412984	20.000	ng	-0.01
39) Acenaphthene-d10	14.469	164	267918	20.000	ng	-0.01
64) Phenanthrene-d10	17.222	188	600837	20.000	ng	#-0.02
76) Chrysene-d12	21.316	240	774310	20.000	ng	-0.02
86) Perylene-d12	23.716	264	911776	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.417	112	751518	106.461	ng	-0.01
7) Phenol-d6	7.011	99	901332	95.978	ng	-0.01
23) Nitrobenzene-d5	8.987	82	559499	72.499	ng	-0.02
42) 2,4,6-Tribromophenol	15.963	330	470964	111.726	ng	-0.01
45) 2-Fluorobiphenyl	13.093	172	1373232	71.897	ng	-0.01
79) Terphenyl-d14	19.792	244	2615421	68.491	ng	-0.01
Target Compounds						
22) Acetophenone	8.652	105	52917	5.037	ng	# 94

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

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