

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112921\
 Data File : BP008140.D
 Acq On : 29 Nov 2021 14:16
 Operator : CG/JU
 Sample : M4844-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 K-1

Quant Time: Nov 29 15:11:45 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 25 01:13:46 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	83325	20.000	ng	-0.01
21) Naphthalene-d8	10.599	136	316225	20.000	ng	-0.01
39) Acenaphthene-d10	14.445	164	206995	20.000	ng	-0.01
64) Phenanthrene-d10	17.204	188	456526	20.000	ng	-0.01
76) Chrysene-d12	21.304	240	508914	20.000	ng	-0.01
86) Perylene-d12	23.698	264	547005	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	596513	115.265	ng	-0.01
7) Phenol-d6	6.987	99	783176	112.105	ng	-0.01
23) Nitrobenzene-d5	8.963	82	530524	76.289	ng	-0.01
42) 2,4,6-Tribromophenol	15.945	330	328916	124.319	ng	-0.01
45) 2-Fluorobiphenyl	13.069	172	1102267	77.349	ng	-0.01
79) Terphenyl-d14	19.780	244	1822198	74.831	ng	0.00
Target Compounds						
75) Fluoranthene	19.228	202	110923	3.416	ng	Qvalue 99
88) Benzo(b)fluoranthene	22.974	252	81256	2.464	ng	# 97
90) Benzo(a)pyrene	23.598	252	57991	2.058	ng	# 95

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

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