

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121721\
 Data File : BP008476.D
 Acq On : 17 Dec 2021 17:53
 Operator : CG/JU
 Sample : M5084-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C2

Quant Time: Dec 18 07:13:17 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP121421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 18 06:49:58 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.728	152	47280	20.000	ng	0.00
21) Naphthalene-d8	10.522	136	175548	20.000	ng	0.00
39) Acenaphthene-d10	14.381	164	119387	20.000	ng	0.00
64) Phenanthrene-d10	17.145	188	260884	20.000	ng	0.00
76) Chrysene-d12	21.257	240	323269	20.000	ng	0.00
86) Perylene-d12	23.621	264	383899	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.328	112	202439	67.618	ng	0.00
7) Phenol-d6	6.934	99	152146	37.839	ng	0.00
23) Nitrobenzene-d5	8.893	82	394325	95.162	ng	0.00
42) 2,4,6-Tribromophenol	15.886	330	271358	154.665	ng	0.00
45) 2-Fluorobiphenyl	12.998	172	850305	90.538	ng	0.00
79) Terphenyl-d14	19.722	244	1472020	79.249	ng	0.00
Target Compounds						
52) Acenaphthene	14.445	154	24216	3.764	ng	99
58) Fluorene	15.439	166	25637	3.022	ng	# 97

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

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