

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092722\
 Data File : BP011813.D
 Acq On : 27 Sep 2022 13:36
 Operator : CG/JU
 Sample : N4809-08
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP3-0923

Quant Time: Sep 28 02:07:12 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 28 02:03:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.910	152	405532	20.000	ng	0.00
21) Naphthalene-d8	10.722	136	1692726	20.000	ng	0.00
39) Acenaphthene-d10	14.557	164	1051499	20.000	ng	0.00
64) Phenanthrene-d10	17.310	188	2165736	20.000	ng	0.00
76) Chrysene-d12	21.392	240	2847453	20.000	ng	0.00
86) Perylene-d12	23.839	264	2965356	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.475	112	3704122	160.265	ng	0.00
7) Phenol-d6	7.081	99	4480551	131.109	ng	0.01
23) Nitrobenzene-d5	9.081	82	2960665	91.546	ng	0.00
42) 2,4,6-Tribromophenol	16.051	330	1484205	209.626	ng	0.00
45) 2-Fluorobiphenyl	13.181	172	6161667	92.668	ng	0.00
79) Terphenyl-d14	19.868	244	11542691	93.208	ng	0.00
Target Compounds						
32) Benzoic acid	10.004	122	40706	8.664	ng	Qvalue 96
84) Bis(2-ethylhexyl)phtha...	21.298	149	29796	2.108	ng	# 99

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

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