

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031522\
 Data File : BP009408.D
 Acq On : 15 Mar 2022 20:46
 Operator : CG/JU
 Sample : N1886-02DL 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 2822-SDL

Quant Time: Mar 16 01:58:22 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 05 01:55:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	311354	20.000	ng	0.00
21) Naphthalene-d8	10.610	136	1232884	20.000	ng	0.00
39) Acenaphthene-d10	14.457	164	779924	20.000	ng	0.00
64) Phenanthrene-d10	17.216	188	1611247	20.000	ng	0.00
76) Chrysene-d12	21.333	240	1708292	20.000	ng	0.00
86) Perylene-d12	23.757	264	1915411	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	232035	10.878	ng	0.01
7) Phenol-d6	7.005	99	300895	11.101	ng	0.02
23) Nitrobenzene-d5	8.969	82	170812	7.631	ng	0.00
42) 2,4,6-Tribromophenol	15.951	330	158049	16.782	ng	0.00
45) 2-Fluorobiphenyl	13.075	172	469716	8.282	ng	-0.01
79) Terphenyl-d14	19.792	244	794416	8.776	ng	0.00
Target Compounds						
74) Di-n-butylphthalate	18.186	149	644453	6.933	ng	# 96
84) Bis(2-ethylhexyl)phtha...	21.251	149	2926379	43.951	ng	# 96

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

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