

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP113020\
 Data File : BP004119.D
 Acq On : 30 Nov 2020 19:13
 Operator : CG/JU
 Sample : L4899-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-1-2

Quant Time: Dec 01 00:25:03 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 23 16:42:41 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.263	152	155959	20.00	ng	0.00
21) Naphthalene-d8	11.092	136	603766	20.00	ng	# 0.00
39) Acenaphthene-d10	14.892	164	372074	20.00	ng	0.00
64) Phenanthrene-d10	17.627	188	816889	20.00	ng	# 0.00
76) Chrysene-d12	21.662	240	770621	20.00	ng	0.00
86) Perylene-d12	24.133	264	655226	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.804	112	1145201	122.55	ng	0.00
7) Phenol-d6	7.457	99	1341770	100.03	ng	0.00
23) Nitrobenzene-d5	9.434	82	1057787	85.80	ng	0.00
42) 2,4,6-Tribromophenol	16.380	330	673018	144.62	ng	0.00
45) 2-Fluorobiphenyl	13.510	172	2307279	86.69	ng	0.00
79) Terphenyl-d14	20.127	244	3330001	83.48	ng	0.00
Target Compounds						
10) Phenol	7.481	94	28192	2.18	ng	# 65
50) Dimethylphthalate	14.316	163	103483	3.56	ng	99

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

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