

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112420\
 Data File : BP004072.D
 Acq On : 24 Nov 2020 12:45
 Operator : CG/JU
 Sample : PB133169BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB133169BL

Quant Time: Nov 24 13:13:39 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	102970	20.00	ng	0.00
21) Naphthalene-d8	11.093	136	388612	20.00	ng	# 0.00
39) Acenaphthene-d10	14.887	164	233848	20.00	ng	0.00
64) Phenanthrene-d10	17.622	188	514732	20.00	ng	0.00
76) Chrysene-d12	21.657	240	508488	20.00	ng	0.00
86) Perylene-d12	24.127	264	325630	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.799	112	851619	138.03	ng	0.00
7) Phenol-d6	7.452	99	1074343	121.31	ng	0.00
23) Nitrobenzene-d5	9.428	82	680510	85.76	ng	0.00
42) 2,4,6-Tribromophenol	16.375	330	385159	131.69	ng	0.00
45) 2-Fluorobiphenyl	13.510	172	1455540	87.02	ng	0.00
79) Terphenyl-d14	20.122	244	2207887	83.89	ng	0.00

Target Compounds	Qvalue

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

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