

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111424\
 Data File : BF140368.D
 Acq On : 14 Nov 2024 17:59
 Operator : RC/JU
 Sample : P4722-10
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-2(0-6)

Quant Time: Nov 15 00:27:19 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF111324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 13 14:40:06 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	124042	20.000	ng	-0.01
21) Naphthalene-d8	8.151	136	464909	20.000	ng	0.00
39) Acenaphthene-d10	9.904	164	252110	20.000	ng	-0.01
64) Phenanthrene-d10	11.392	188	459695	20.000	ng	-0.01
76) Chrysene-d12	14.045	240	260161	20.000	ng	0.00
86) Perylene-d12	15.545	264	261026	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	541049	74.473	ng	0.00
7) Phenol-d6	6.493	99	469938	47.744	ng	-0.01
23) Nitrobenzene-d5	7.434	82	850173	95.280	ng	0.00
42) 2,4,6-Tribromophenol	10.698	330	325204	129.716	ng	0.00
45) 2-Fluorobiphenyl	9.228	172	1421774	90.658	ng	0.00
79) Terphenyl-d14	12.980	244	1479681	98.724	ng	0.00
Target Compounds						
59) 2,3,4,6-Tetrachlorophenol	10.233	232	8775	2.222	ng	Qvalue # 97
70) Pentachlorophenol	11.198	266	27258	9.787	ng	96

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

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