

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112921\
 Data File : BP008138.D
 Acq On : 29 Nov 2021 13:08
 Operator : CG/JU
 Sample : M4798-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-1-3.0

Quant Time: Nov 29 13:41:07 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 25 01:13:46 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	96990	20.000	ng	-0.01
21) Naphthalene-d8	10.605	136	335509	20.000	ng	# 0.00
39) Acenaphthene-d10	14.463	164	197467	20.000	ng	0.00
64) Phenanthrene-d10	17.222	188	403398	20.000	ng	# 0.00
76) Chrysene-d12	21.310	240	478494	20.000	ng	0.00
86) Perylene-d12	23.698	264	526117	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	509443	84.571	ng	-0.01
7) Phenol-d6	6.987	99	665545	81.845	ng	-0.01
23) Nitrobenzene-d5	8.963	82	475089	64.391	ng	-0.01
42) 2,4,6-Tribromophenol	15.957	330	242794	96.196	ng	0.00
45) 2-Fluorobiphenyl	13.081	172	785844	57.805	ng	0.00
79) Terphenyl-d14	19.781	244	1297022	56.650	ng	0.00
Target Compounds						
						Qvalue
38) 1-Methylnaphthalene	12.493	142	192634	16.280	ng	# 97
49) Acenaphthylene	14.198	152	65730	3.766	ng	# 1
52) Acenaphthene	14.528	154	102196	9.824	ng	89
55) Dibenzofuran	14.863	168	66627	3.748	ng	# 1
58) Fluorene	15.516	166	80319	5.561	ng	# 61
71) Phenanthrene	17.263	178	83169	3.905	ng	# 63
75) Fluoranthene	19.233	202	80921	2.820	ng	76
78) Pyrene	19.586	202	233238	5.466	ng	# 94

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

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