

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042424\
 Data File : BF137722.D
 Acq On : 24 Apr 2024 17:12
 Operator : CG\JU
 Sample : P2208-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LNG-11

Quant Time: Apr 24 17:49:01 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 24 03:57:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.063	152	263474	20.000	ng	0.00
21) Naphthalene-d8	8.345	136	1028691	20.000	ng	0.00
39) Acenaphthene-d10	10.104	164	541711	20.000	ng	-0.01
64) Phenanthrene-d10	11.604	188	890988	20.000	ng	# 0.00
76) Chrysene-d12	14.251	240	499433	20.000	ng	# 0.00
86) Perylene-d12	15.827	264	544453	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.692	112	1856906	107.058	ng	0.01
7) Phenol-d6	6.675	99	2175516	98.770	ng	0.00
23) Nitrobenzene-d5	7.622	82	1499755	77.391	ng	0.00
42) 2,4,6-Tribromophenol	10.898	330	640489	117.171	ng	0.00
45) 2-Fluorobiphenyl	9.422	172	2773207	76.494	ng	0.00
79) Terphenyl-d14	13.192	244	2876748	92.711	ng	0.00
Target Compounds						
15) Benzyl Alcohol	7.186	79	45273	2.748	ng	98
20) 3+4-Methylphenols	7.439	107	56614	2.725	ng	93

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

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