

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031325\
 Data File : BF141943.D
 Acq On : 13 Mar 2025 16:14
 Operator : RC/JU
 Sample : Q1539-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TAPIAL3-MW03D-031025-00-T1

Quant Time: Mar 13 16:47:28 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF031025.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 10 15:46:22 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	170628	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	672168	20.000	ng	0.00
39) Acenaphthene-d10	9.910	164	390587	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	678228	20.000	ng	0.00
76) Chrysene-d12	14.027	240	432058	20.000	ng	-0.01
86) Perylene-d12	15.498	264	372025	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.492	112	600091	58.697	ng	0.00
7) Phenol-d6	6.487	99	461285	35.437	ng	-0.02
23) Nitrobenzene-d5	7.434	82	1194711	100.025	ng	-0.01
42) 2,4,6-Tribromophenol	10.704	330	789341	159.298	ng	0.00
45) 2-Fluorobiphenyl	9.233	172	2464310	95.952	ng	0.00
79) Terphenyl-d14	12.980	244	2674361	91.514	ng	0.00
Target Compounds						
						Qvalue
52) Acenaphthene	9.939	154	54270	2.353	ng	# 88
55) Dibenzofuran	10.110	168	85562	2.596	ng	99
58) Fluorene	10.457	166	112667	4.432	ng	99
73) Carbazole	11.621	167	69087	2.180	ng	99

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

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