

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031325\
 Data File : BF141951.D
 Acq On : 13 Mar 2025 20:10
 Operator : RC/JU
 Sample : Q1547-10
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-620-JB-COMP-02

Manual Integrations
 APPROVED

Reviewed By :Anahy Claudio 03/14/2025
 Supervised By :Jagrut Upadhyay 03/14/2025

Quant Time: Mar 14 10:40:04 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	169785	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	672608	20.000	ng	0.00
39) Acenaphthene-d10	9.910	164	391126	20.000	ng	0.00
64) Phenanthrene-d10	11.392	188	701302	20.000	ng	-0.01
76) Chrysene-d12	14.027	240	403778	20.000	ng	-0.01
86) Perylene-d12	15.498	264	366915	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	1455596	143.083	ng	0.02
7) Phenol-d6	6.498	99	1655739	127.831	ng	0.00
23) Nitrobenzene-d5	7.440	82	1302918	109.013	ng	0.00
42) 2,4,6-Tribromophenol	10.698	330	911372	183.672	ng	0.00
45) 2-Fluorobiphenyl	9.234	172	2733900	106.302	ng	0.00
79) Terphenyl-d14	12.980	244	3360141	123.033	ng	0.00
Target Compounds						
32) Benzoic acid	7.851	122	18822m	2.667	ng	Qvalue
50) Dimethylphthalate	9.622	163	59044	2.156	ng	98
77) Benzidine	12.727	184	53665	9.526	ng	99

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

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