

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF032825\
 Data File : BF142163.D
 Acq On : 28 Mar 2025 16:35
 Operator : RC/JU
 Sample : Q1662-04
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-6

Manual Integrations
 APPROVED

Reviewed By :Anahy Claudio 03/31/2025
 Supervised By :Jagrut Upadhyay 03/31/2025

Quant Time: Mar 28 16:58:55 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	6.869	152	141234	20.000	ng	-0.01
21) Naphthalene-d8	8.151	136	583550	20.000	ng	-0.01
39) Acenaphthene-d10	9.904	164	329645	20.000	ng	-0.01
64) Phenanthrene-d10	11.392	188	523823	20.000	ng	-0.01
76) Chrysene-d12	14.033	240	304790	20.000	ng	0.00
86) Perylene-d12	15.510	264	323130	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	447011	52.823	ng	0.00
7) Phenol-d6	6.492	99	518857	48.156	ng	-0.01
23) Nitrobenzene-d5	7.428	82	394210	38.017	ng	-0.02
42) 2,4,6-Tribromophenol	10.692	330	187688	44.880	ng	-0.01
45) 2-Fluorobiphenyl	9.228	172	866233	39.963	ng	-0.01
79) Terphenyl-d14	12.974	244	783457	38.003	ng	-0.01
Target Compounds						
77) Benzidine	12.733	184	23926m	5.626	ng	Qvalue

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

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