

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110722\
 Data File : BP012531.D
 Acq On : 07 Nov 2022 15:00
 Operator : CG/JU
 Sample : N5353-11
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBVV1

Quant Time: Nov 07 21:41:10 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 04 02:53:36 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	268507	20.000	ng/ul	0.00
20) Naphthalene-d8	10.575	136	1325759	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.428	164	921243	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.192	188	2088529	20.000	ng/ul	0.00
79) Chrysene-d12	21.286	240	1844135	20.000	ng/ul	0.00
88) Perylene-d12	23.668	264	1678981	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	31351	4.701	ng/uL	0.00
4) Pyridine-d5	3.658	84	142299	7.428	ng/ul	0.00
7) Phenol-d5	6.958	99	141020	5.856	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.116	67	414708	28.846	ng/ul	0.00
11) 2-Chlorophenol-d4	7.310	132	410716	23.353	ng/ul	0.00
15) 4-Methylphenol-d8	8.499	113	301189	15.813	ng/ul	0.00
21) Nitrobenzene-d5	8.946	128	298137	34.678	ng/ul	0.00
24) 2-Nitrophenol-d4	9.669	143	314702	35.267	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	575707	29.770	ng/ul	0.00
31) 4-Chloroaniline-d4	10.728	131	723973	25.143	ng/ul	0.00
46) Dimethylphthalate-d6	13.845	166	2205121	34.966	ng/ul	0.00
49) Acenaphthylene-d8	14.122	160	2343693	32.520	ng/ul	0.00
54) 4-Nitrophenol-d4	14.628	143	217	0.018	ng/ul	-0.03
60) Fluorene-d10	15.428	176	1944086	36.006	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	98486	9.057	ng/ul	0.00
73) Anthracene-d10	17.292	188	3295631	36.622	ng/ul	0.00
81) Pyrene-d10	19.533	212	3795377	40.992	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.515	264	3032363	36.963	ng/ul	0.00

Target Compounds Qvalue

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

