

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009762.D
 Acq On : 11 Apr 2022 21:34
 Operator : CG/JU
 Sample : N2338-20
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J86

Quant Time: Apr 12 01:53:36 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 07 19:16:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	108105	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.781	136	462357	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.604	164	324983	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.351	188	735760	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.433	240	791343	20.000	ng/u1	#-0.01
88) Perylene-d12	23.904	264	788169	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	15077	6.224	ng/uL	0.00
4) Pyridine-d5	3.805	84	74640	11.176	ng/u1	0.00
7) Phenol-d5	7.122	99	78994	8.323	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	219774	38.896	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.498	132	215117	30.048	ng/u1	0.00
15) 4-Methylphenol-d8	8.669	113	149971	18.929	ng/u1	-0.01
21) Nitrobenzene-d5	9.128	128	139912	41.963	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.851	143	146437	40.060	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.392	165	253780	33.177	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.904	131	196559	18.042	ng/u1	-0.01
46) Dimethylphthalate-d6	14.010	166	1043424	40.874	ng/u1	-0.01
49) Acenaphthylene-d8	14.292	160	1121546	37.084	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.775	143	41476	8.959	ng/u1	-0.01
60) Fluorene-d10	15.592	176	897463	41.791	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.698	200	178720	41.003	ng/u1	-0.01
73) Anthracene-d10	17.451	188	1565827	44.449	ng/u1	-0.01
81) Pyrene-d10	19.680	212	1936022	44.795	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.745	264	1907130	46.464	ng/u1	-0.02

Target Compounds Qvalue

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

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