

Data Path : Z:\HPCHEM1\BNA E\DATA\BE112215\
 Data File : BE091292.D
 Acq On : 22 Nov 2015 10:12
 Operator : UM/NP
 Sample : PB86665BL
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SBLK65

Manual Integrations
 APPROVED

Mmdadoda
 11/23/2015 2:14:40 PM

Quant Time: Nov 23 00:49:15 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM02.2-EPA-SIM-BE102715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 23 00:41:30 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	2985	0.40	ng/ul	-0.03
4) Naphthalene-d8	9.54	136	14790	0.40	ng/ul	-0.02
8) Acenaphthene-d10	13.36	164	7961	0.40	ng/ul	0.00
12) Phenanthrene-d10	16.08	188	22101m	0.40	ng/ul	-0.01
18) Chrysene-d12	20.22	240	43223	0.40	ng/ul	0.00
22) Perylene-d12	22.11	264	29662m	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	2.91	96	25119	11.70	ng/uL	-0.04
6) 2-Methylnaphthalene-d10	11.04	152	5457	0.27	ng/ul	-0.01
16) Fluoranthene-d10	18.07	212	13775	0.22	ng/ul	0.00
Target Compounds						
13) Pentachlorophenol	15.82	266	1327	0.63	ng/ul	Ovalue 95
17) Fluoranthene	18.09	202	78631	0.26	ng/ul	88
19) Pyrene	18.47	202	94985	0.19	ng/ul#	76
20) Benzo(a)anthracene	20.20	228	14620	0.13	ng/ul	96
21) Chrysene	20.26	228	19579	0.15	ng/ul	97
23) Benzo(b)fluoranthene	21.56	252	6064	0.07	ng/ul	97
24) Benzo(k)fluoranthene	21.58	252	6202m	0.07	ng/ul	
25) Benzo(a)pyrene	22.01	252	4438	0.05	ng/ul	96

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

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